

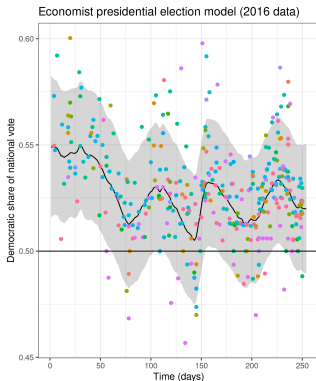
Approximate data deletion and replication with the Bayesian influence function

Ryan Giordano (rgiordano@berkeley.edu, UC Berkeley), Tamara Broderick (MIT)

April 2024

Theory and Foundations of Statistics in the Era of Big Data

Economist 2016 Election Model [Gelman and Heidemanns, 2020]



A time series model to predict the 2016 US presidential election outcome from polling data.

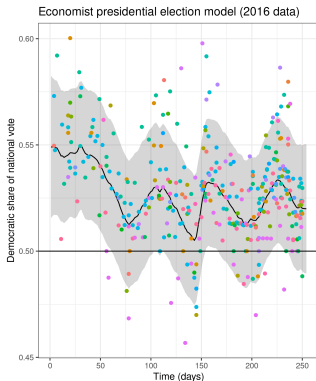
Model:

- $X = x_1, \dots, x_N =$ Polling data ($N = 361$).
- $\theta =$ Lots of random effects (day, pollster, etc.)
- $f(\theta) =$ Democratic % of vote on election day

Typically, we compute Markov chain Monte Carlo (MCMC) draws from the posterior $p(\theta|X)$.

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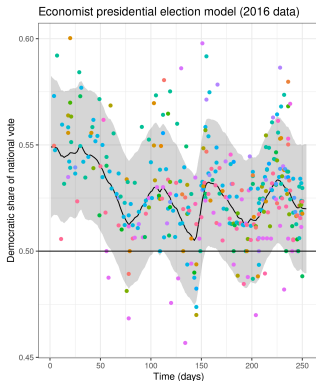
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Re-fit with data points removed one at a time

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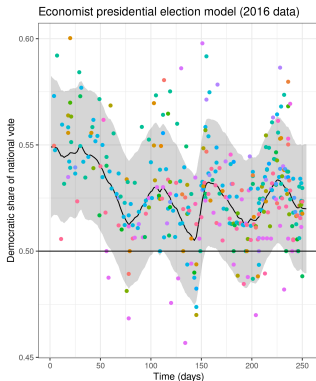
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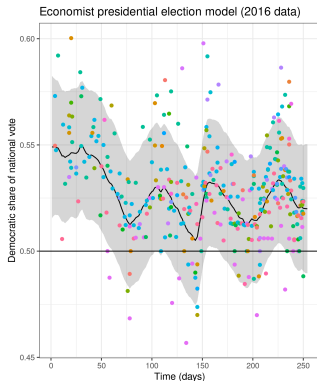
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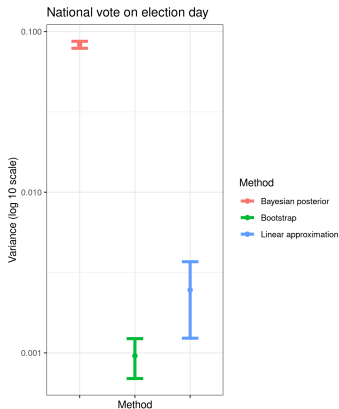
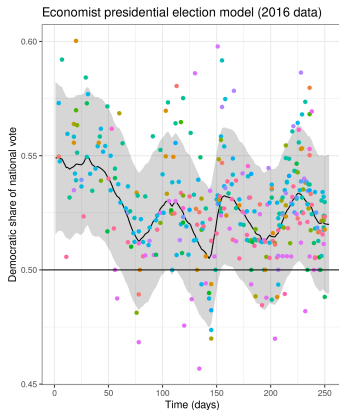
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Problem: Each MCMC run takes about 10 hours (Stan, six cores).

We propose: Use posterior draws based on the full data, to form a linear approximation to *data reweightings*.

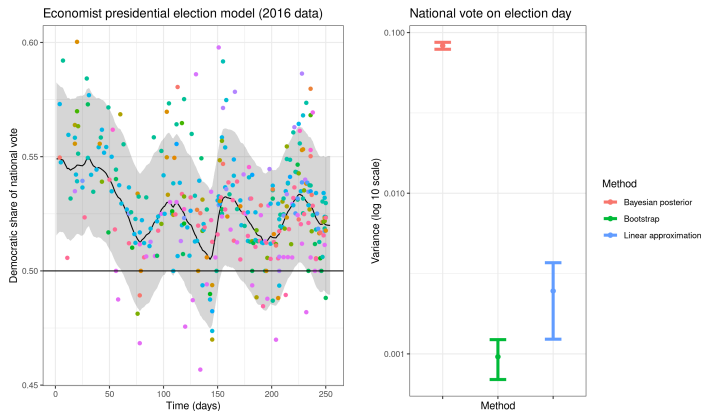
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Compute time for 100 bootstraps: 51 days

Compute time for the linear approximation: Seconds
(But note the approximation has some error)

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- A trick question, and some implications of different weightings.

Data re-weighting.

Augment the problem with *data weights* w_1, \dots, w_N . We can write $\mathbb{E}_{p(\theta|X, w)} [f(\theta)]$.

$$\ell_n(\theta) := \log p(x_n|\theta)$$

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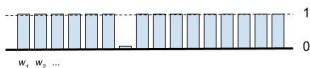
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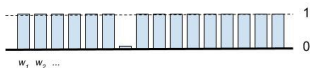
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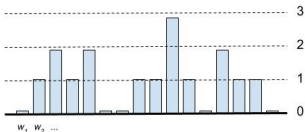
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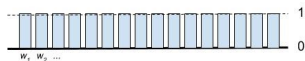
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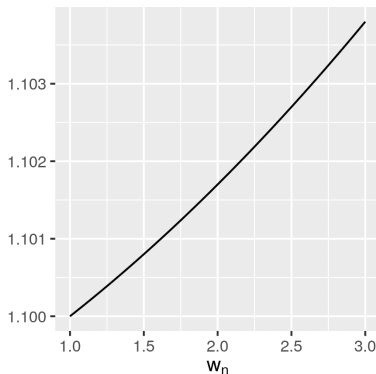
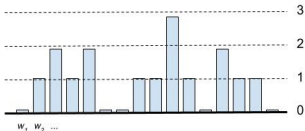
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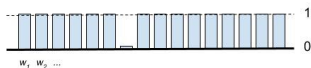
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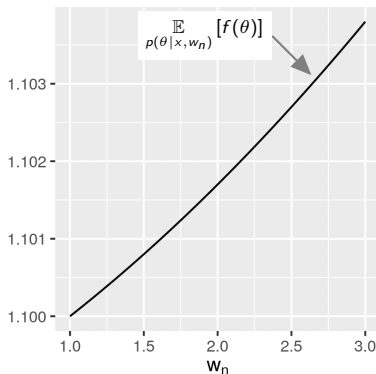
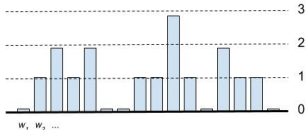
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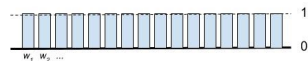
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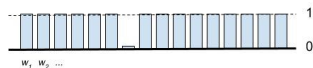
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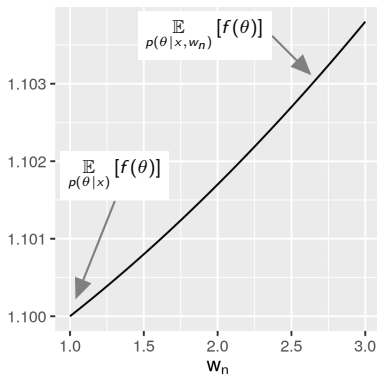
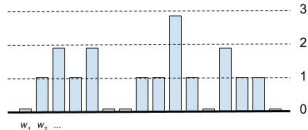
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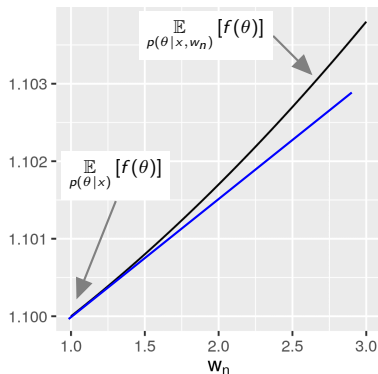
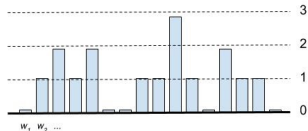
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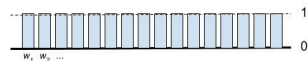
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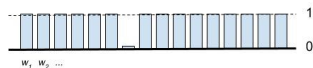
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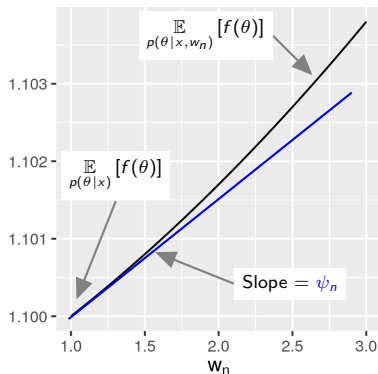
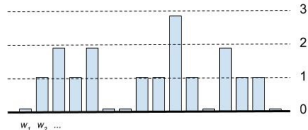
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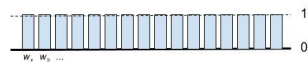
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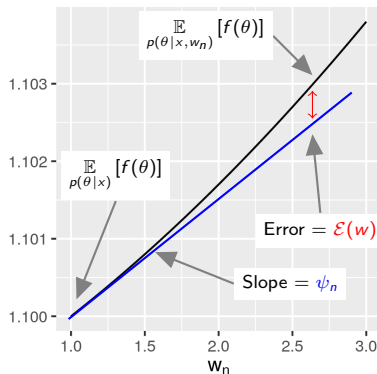
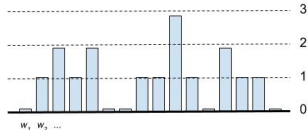
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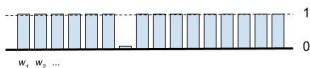
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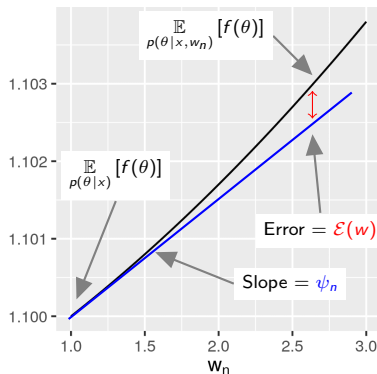
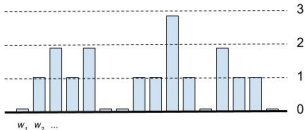
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The re-scaled slope $N\psi_n$ is known as the “influence function” at data point x_n .

$$\mathbb{E}_{p(\theta|X,w)} [f(\theta)] - \mathbb{E}_{p(\theta|X)} [f(\theta)] = \sum_{n=1}^N \psi_n (w_n - 1) + \mathcal{E}(w)$$

How can we use the approximation?

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Influential subsets: Approximate maximum influence perturbation (AMIP).

Let $W_{(-K)}$ denote weights leaving out K points.

$$\max_{w \in W_{(-K)}} \left(\mathbb{E}_{p(\theta|x,w)}[f(\theta)] - \mathbb{E}_{p(\theta|x)}[f(\theta)] \right) \approx - \sum_{n=1}^K \psi_{(n)}.$$

Expressions for the slope and error

How to compute the slopes ψ_n ? How large is the error $\mathcal{E}(w)$?

For simplicity, for the remainder of the presentation, we will consider a single weight.

$$\mathbb{E}_{p(\theta|X, w_n)} [f(\theta)] - \mathbb{E}_{p(\theta|X)} [f(\theta)] = \psi_n (w_n - 1) + \mathcal{E}(w_n)$$

Let an overbar mean posterior-mean zero (e.g., $\bar{f}(\theta) := f(\theta) - \mathbb{E}_{p(\theta|X)} [f(\theta)]$).

By dominated convergence and the mean value theorem, for some $\tilde{w}_n \in [0, w_n]$:

$$\begin{aligned} \psi_n &= \underbrace{\mathbb{E}_{p(\theta|X)} [\bar{f}(\theta) \bar{\ell}_n(\theta)]}_{\text{Estimatable with MCMC!}} \\ &= O_p(N^{-1}) \text{ under a BCLT} \end{aligned} \quad \begin{aligned} \mathcal{E}(w_n) &= \frac{1}{2} \underbrace{\mathbb{E}_{p(\theta|X, \tilde{w}_n)} [\bar{f}(\theta) \bar{\ell}_n(\theta) \bar{\ell}_n(\theta)]}_{\text{Cannot compute directly (don't know } \tilde{w})}} (w_n - 1)^2 \\ &= O_p(N^{-2}) \text{ under a BCLT} \end{aligned}$$

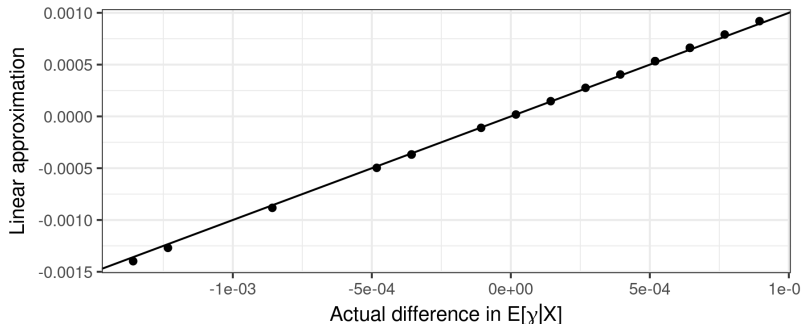
\Rightarrow The map $w_n \mapsto N \left(\mathbb{E}_{p(\theta|X, w_n)} [f(\theta)] - \mathbb{E}_{p(\theta|X)} [f(\theta)] \right)$ becomes linear as $N \rightarrow \infty$.

(See [Kass et al., 1990] for a *particular weight*, [?] for a kind of uniform convergence over datapoints.)

Low dimensional problems

Example: **Negative binomial models with an unknown parameter γ .**

Negative Binomial model
leaving out single datapoints with $N = 800$



The map $w_n \mapsto N \left(\mathbb{E}_{p(\gamma|X, w_n)} [\gamma] - \mathbb{E}_{p(\gamma|X)} [\gamma] \right)$ becomes linear as $N \rightarrow \infty$.

High dimensional problems

What about when the posterior doesn't obey a BCLT?

Suppose that $p(\lambda|X)$ does not concentrate.

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Can we save the approximation when *some* parameters concentrate?

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If the observations per random effect remains bounded as $N \rightarrow \infty$, then

- Parameter λ (“local”) grows in dimension with N .
- Parameter γ (“global”) is finite-dimensional.
- Marginally $p(\lambda|X)$ does not concentrate.
- Marginally, $p(\gamma|X)$ concentrates.

Example: **Poisson model with random effects (REs) λ and fixed effects γ .**

If the observations per random effect remains bounded as $N \rightarrow \infty$, then

- Parameter λ (“local”) grows in dimension with N .
- Parameter γ (“global”) is finite-dimensional.
- Marginally $p(\lambda|X)$ does not concentrate.
- Marginally, $p(\gamma|X)$ concentrates.

Can we save the approximation when *some* parameters concentrate?

\Rightarrow Does the residual vanish asymptotically for $w_n \mapsto \mathbb{E}_{p(\gamma|X, w_n)}[\gamma]$?

High dimensional problems

We assume that $p(\gamma|X)$ concentrates but $p(\lambda|X)$ does not. By our series expansion:

$$\begin{aligned} & \mathbb{E}_{p(\gamma, \lambda|X, w_n)} [\gamma] - \mathbb{E}_{p(\gamma, \lambda|X)} [\gamma] = \\ & \quad \psi_n(w_n - 1) + \mathcal{E}(w_n) \\ = & \mathbb{E}_{p(\gamma, \lambda|X)} [\tilde{\gamma} \bar{\ell}_n(\gamma, \lambda)] (w_n - 1) + \frac{1}{2} \mathbb{E}_{p(\gamma, \lambda|X, \tilde{w}_n)} [\tilde{\gamma} \bar{\ell}_n(\gamma, \lambda)^2] (w_n - 1)^2 \end{aligned}$$

$$\psi_n = O_p(N^{-1})$$

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High dimensional problems

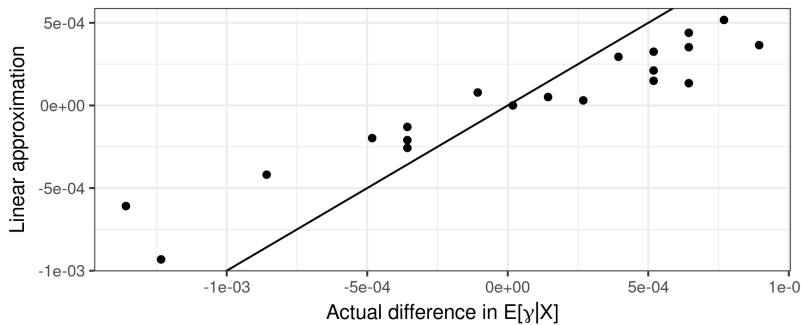
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The map $w_n \mapsto N \left(\mathbb{E}_{p(\gamma|X, w_n)} [\gamma] - \mathbb{E}_{p(\gamma|X)} [\gamma] \right)$ remains non-linear as $N \rightarrow \infty$.

Experiments

Poisson random effect model
leaving out single datapoints with N = 800



A contradiction?

Negative binomial observations.

Asymptotically linear in w .

Poisson observations with random effects.

Asymptotically non-linear in w .

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With a constant regressor, Gamma REs, and one RE per observation, these are the same model, with the same $p(\gamma|X)$.

Is $\mathbb{E}_{p(\gamma|X,w)}[\gamma]$ linear in the data weights or not?

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$$\log p(X|\gamma, w^m) = \sum_{n=1}^N w_n^m \log p(x_n|\gamma)$$

Poisson observations with random effects.

Asymptotically non-linear in w .

$$\log p(X|\gamma, \lambda, w^c) = \sum_{n=1}^N w_n^c \log p(x_n|\lambda, \gamma)$$

With a constant regressor, Gamma REs, and one RE per observation, these are the same model, with the same $p(\gamma|X)$.

Is $\mathbb{E}_{p(\gamma|X, w)}[\gamma]$ linear in the **data weights** or not?

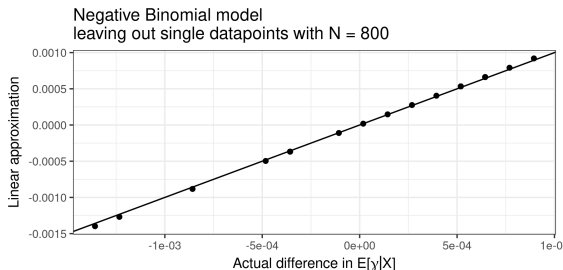
Trick question! We weight a log likelihood contribution, not a datapoint.

The two weightings are not equivalent in general.

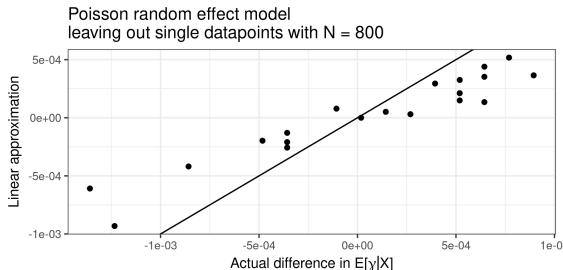
Experimental results

Our results were actually computed on **identical datasets** with $G = N$ and $g_n = n$.

Approximation based
on $\log p(x_n|\gamma)$.



Approximation based
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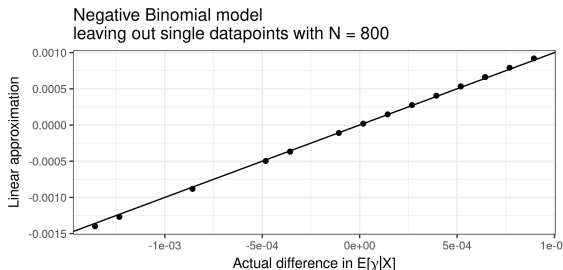


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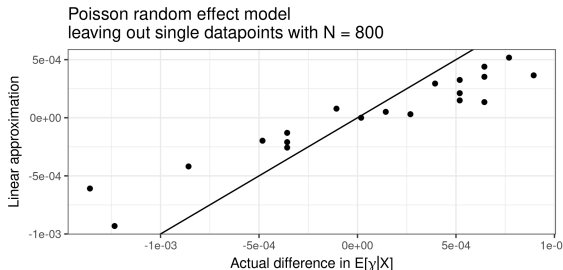
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 $\gamma, \lambda \sim p(\gamma, \lambda|X)$
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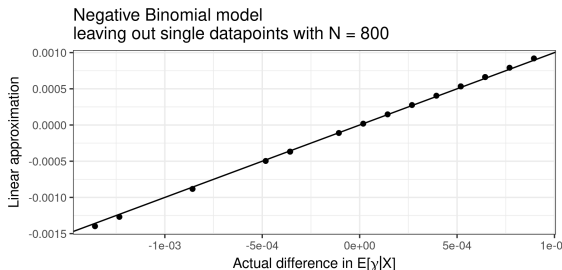


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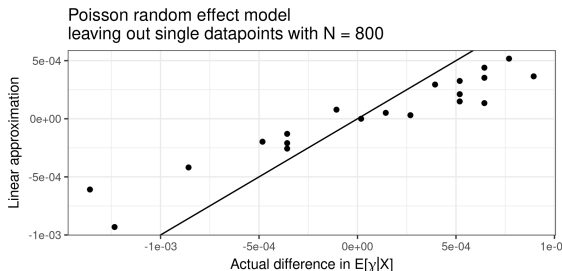
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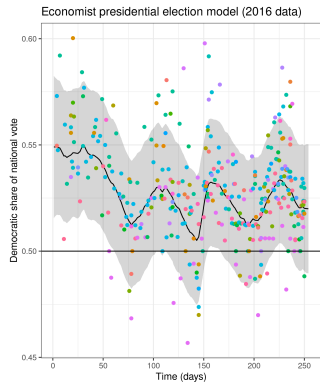
Approximation based
on $\log p(x_n|\gamma, \lambda)$.

Computable from
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May still be useful
when $p(\lambda|X)$ is *some-
what* concentrated.

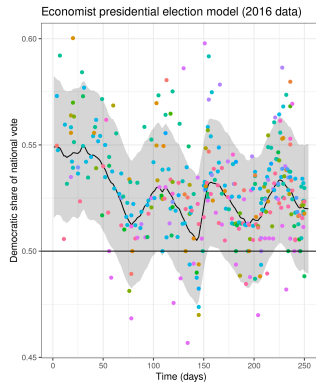


Observations and consequences



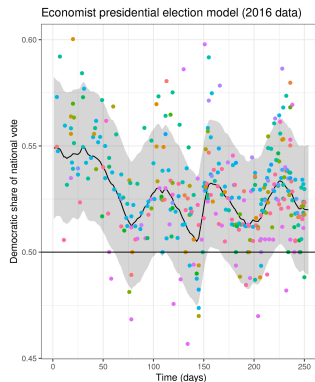
Observations and consequences

- When $\log p(x_n|\gamma, \lambda)$ is the exchangeable unit, our results are problematic for
 - Linear approximations (IJ, AMIP, approx. CV)
 - The nonparametric bootstrap
 - All of the above for Bayes-like optimization procedures (VB, the EM algorithm)



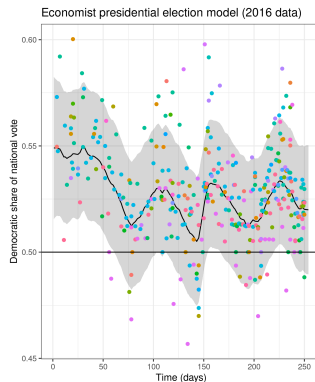
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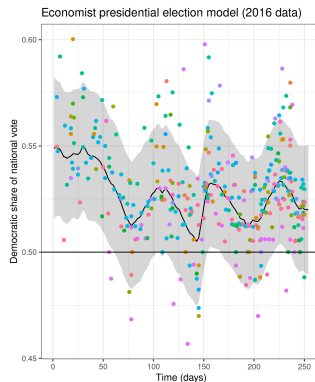
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- There may be multiple ways to define “exchangeable unit” in a given problem.
- But without nesting, $\log p(x_n|\gamma, \lambda)$ may be the natural model-free exchangeable unit.



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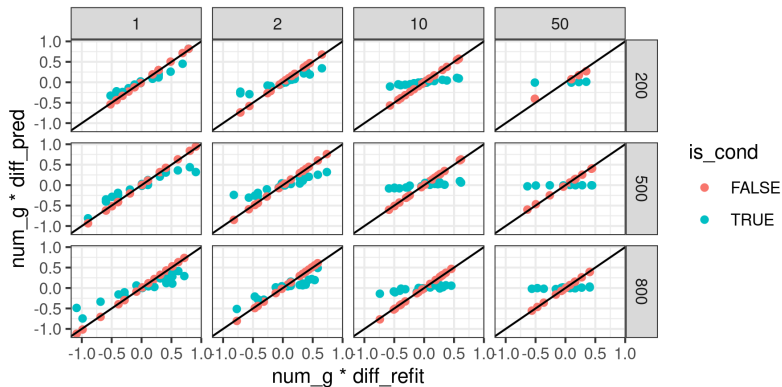
Supplemental slides

Non-equivalence of weighting (nonlinearity of marginalization)

Consider a single datapoint.

$$\begin{aligned}\log p(x_n|\gamma, w_c) &= \\ \log \left(\int p(x_n|\gamma, \lambda, w_c) p(\lambda|\gamma) d\lambda \right) &= \\ \log \left(\int p(x_n|\gamma, \lambda)^{w_c} p(\lambda|\gamma) d\lambda \right) &\neq \\ \log \left(\int p(x_n|\gamma, \lambda) p(\lambda|\gamma) d\lambda \right)^{w_c} &= \\ w_c \log p(\lambda|\gamma)\end{aligned}$$

Extended experimental results



Extended experimental results

